

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111821\
 Data File : BP008022.D
 Acq On : 18 Nov 2021 18:27
 Operator : CG/JU
 Sample : M4666-01RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CONCRETERE

Quant Time: Nov 19 02:24:12 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 15 05:36:49 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	100390	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	331533	20.000	ng	#-0.01
39) Acenaphthene-d10	14.469	164	210017	20.000	ng	-0.01
64) Phenanthrene-d10	17.228	188	463313	20.000	ng	#-0.01
76) Chrysene-d12	21.322	240	571384	20.000	ng	#-0.01
86) Perylene-d12	23.721	264	646239	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	44486	7.189	ng	0.00
7) Phenol-d6	7.011	99	268001	32.553	ng	-0.01
23) Nitrobenzene-d5	8.987	82	340836	55.015	ng	-0.02
42) 2,4,6-Tribromophenol	15.969	330	4752	1.438	ng	0.00
45) 2-Fluorobiphenyl	13.093	172	752623	50.268	ng	-0.01
79) Terphenyl-d14	19.792	244	1367104	48.515	ng	-0.01
Target Compounds						
						Qvalue
80) Butylbenzylphthalate	20.475	149	41089	2.912	ng	92
81) Benzo(a)anthracene	21.304	228	96209	2.558	ng	99
83) Chrysene	21.357	228	114082	3.164	ng	98
84) Bis(2-ethylhexyl)phtha...	21.239	149	52501	2.606	ng	# 95
85) Di-n-octyl phthalate	22.163	149	130197	3.507	ng	# 99
88) Benzo(b)fluoranthene	22.992	252	126653	3.292	ng	100
89) Benzo(k)fluoranthene	23.039	252	125868	3.275	ng	96
90) Benzo(a)pyrene	23.616	252	74907	2.277	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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